

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121018\
 Data File : BN003777.D
 Acq On : 10 Dec 2018 14:11
 Operator : JU/SJ
 Sample : J6155-23
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4132

Quant Time: Dec 11 03:57:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 11 03:42:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4373	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	17795	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	9324	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	20349	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	16501	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	14045	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	10092	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	20417	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.68	128	42940	0.859	ng/ul	100
7) Acenaphthylene	14.19	152	34364	0.818	ng/ul	100
8) Acenaphthene	14.53	153	28156	0.765	ng/ul	99
9) Fluorene	15.52	166	37427	0.873	ng/ul	100
11) Pentachlorophenol	16.86	266	9049	2.469	ng/ul	100
12) Phenanthrene	17.25	178	49725	0.745	ng/ul	99
13) Anthracene	17.35	178	26263	0.478	ng/ul	100
17) Pyrene	19.63	202	110810	1.872	ng/ul	98
18) Benzo(a)anthracene	21.38	228	41051	0.788	ng/ul	99
19) Chrysene	21.43	228	61198	1.026	ng/ul	100
22) Benzo(k)fluoranthene	23.06	252	26909	0.475	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	38914	0.674	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.06	278	1436	0.031	ng/ul#	53
26) Benzo(g,h,i)perylene	26.78	276	43752	0.873	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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